

11/539,485

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PASSWORD:

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NEWS	4	APR 04 STN AnaVist \$500 visualization usage credit offered
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NEWS	6	MAY 11 KOREAPAT updates resume
NEWS	7	MAY 19 Derwent World Patents Index to be reloaded and enhanced
NEWS	8	MAY 30 IPC 8 Rolled-up Core codes added to CA/CAPLUS and USPATFULL/USPAT2
NEWS	9	MAY 30 The F-Term thesaurus is now available in CA/CAPLUS
NEWS	10	JUN 02 The first reclassification of IPC codes now complete in INPADOC
NEWS	11	JUN 26 TULSA/TULSA2 reloaded and enhanced with new search and and display fields
NEWS	12	JUN 28 Price changes in full-text patent databases EPFULL and PCTFULL
NEWS	13	JUL 11 CHEMSAFE reloaded and enhanced
NEWS	14	JUL 14 FSTA enhanced with Japanese patents
NEWS	15	JUL 19 Coverage of Research Disclosure reinstated in DWPI
NEWS	16	AUG 09 INSPEC enhanced with 1898-1968 archive
NEWS EXPRESS		JUNE 30 CURRENT WINDOWS VERSION IS V8.01b, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 26 JUNE 2006.
NEWS HOURS		STN Operating Hours Plus Help Desk Availability
NEWS LOGIN		Welcome Banner and News Items
NEWS IPC8		For general information regarding STN implementation of IPC 8
NEWS X25		X.25 communication option no longer available

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 14:23:04 ON 14 AUG 2006

=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION

11/539,485

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 14:23:13 ON 14 AUG 2006
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STRUCTURE FILE UPDATES: 13 AUG 2006 HIGHEST RN 901009-82-3
DICTIONARY FILE UPDATES: 13 AUG 2006 HIGHEST RN 901009-82-3

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TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

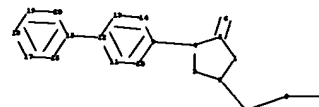
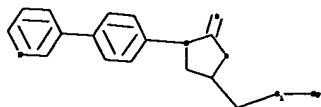
Please note that search-term pricing does apply when
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REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10539485-1.str



chain nodes :

6 7 8 22

ring nodes :

1 2 3 4 5 9 10 11 12 13 14 15 16 17 18 19 20

chain bonds :

11/539,485

2-7 4-9 5-6 7-8 8-22 12-15

ring bonds :

1-2 1-5 2-3 3-4 4-5 9-10 9-14 10-11 11-12 12-13 13-14 15-16 15-20
16-17 17-18 18-19 19-20

exact/norm bonds :

1-2 1-5 2-3 3-4 4-5 4-9 5-6 7-8 8-22

exact bonds :

2-7 12-15

normalized bonds :

9-10 9-14 10-11 11-12 12-13 13-14 15-16 15-20 16-17 17-18 18-19 19-20

isolated ring systems :

containing 9 :

G1:O,S

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:CLASS 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom 22:Atom

Generic attributes :

22:

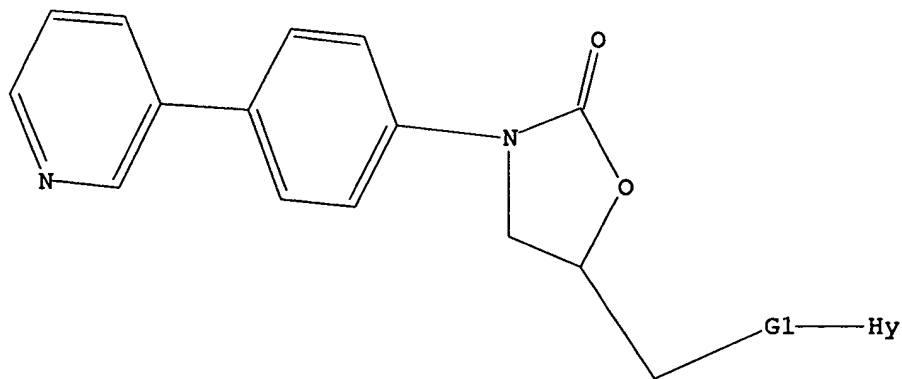
Saturation : Unsaturated

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



G1 O,S

Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss sam

SAMPLE SEARCH INITIATED 14:23:36 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 12 TO ITERATE

100.0% PROCESSED 12 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

11/539,485

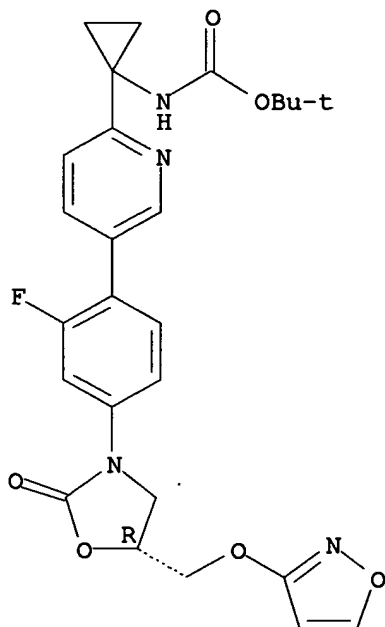
FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 33 TO 447
PROJECTED ANSWERS: 1 TO 80

L2 1 SEA SSS SAM L1

=> d scan

L2 1 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Carbamic acid, [1-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]cyclopropyl]-, 1,1-dimethylethyl ester (9CI)
MF C26 H27 F N4 O6

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> s l1 sss ful
FULL SEARCH INITIATED 14:23:51 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 297 TO ITERATE

100.0% PROCESSED 297 ITERATIONS 12 ANSWERS
SEARCH TIME: 00.00.01

L3 12 SEA SSS FUL L1

=> file caplus
COST IN U.S. DOLLARS SINCE FILE TOTAL

	ENTRY	SESSION
FULL ESTIMATED COST	166.94	167.15

FILE 'CAPLUS' ENTERED AT 14:23:58 ON 14 AUG 2006
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FILE COVERS 1907 - 14 Aug 2006 VOL 145 ISS 8
 FILE LAST UPDATED: 13 Aug 2006 (20060813/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/infopolicy.html>

=> s 13

L4 3 L3

=> d 14 1-3 bib hitstr

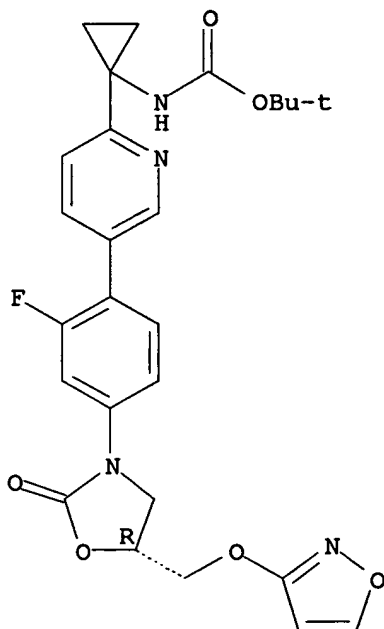
L4 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN
 AN 2005:58198 CAPLUS
 DN 142:155938
 TI Preparation of cyclopropyl group substituted oxazolidinones as antibiotics
 IN Fukuda, Yasumichi
 PA Merck & Co., Inc., USA; Kyorin Pharmaceutical Co., Ltd.
 SO PCT Int. Appl., 85 pp.
 CODEN: PIXXD2
 DT Patent
 LA English
 FAN.CNT 2

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2005005420	A1	20050120	WO 2004-US20737	20040629
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
	RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
	AU 2004256085	A1	20050120	AU 2004-256085	20040629
	CA 2529293	AA	20050120	CA 2004-2529293	20040629
	EP 1654259	A1	20060510	EP 2004-777199	20040629
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,				

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IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK
CN 1816545 A 20060809 CN 2004-80018905 20040629
PRAI US 2003-483904P P 20030702
US 2004-546980P P 20040224
WO 2004-US20737 W 20040629
OS CASREACT 142:155938; MARPAT 142:155938
IT 827627-86-1P
RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(antibacterial agent; preparation of cyclopropyl-oxazolidinones as antibiotics)
RN 827627-86-1 CAPLUS
CN Carbamic acid, [1-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]cyclopropyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

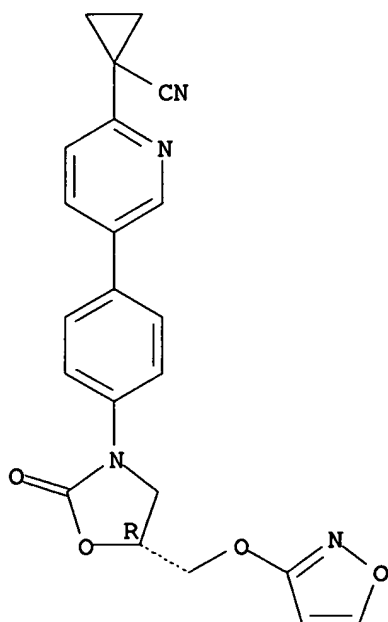
Absolute stereochemistry.



IT 827627-81-6P 827627-87-2P 827627-90-7P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(antibacterial agent; preparation of cyclopropyl-oxazolidinones as antibiotics)
RN 827627-81-6 CAPLUS
CN Cyclopropanecarbonitrile, 1-[5-[4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

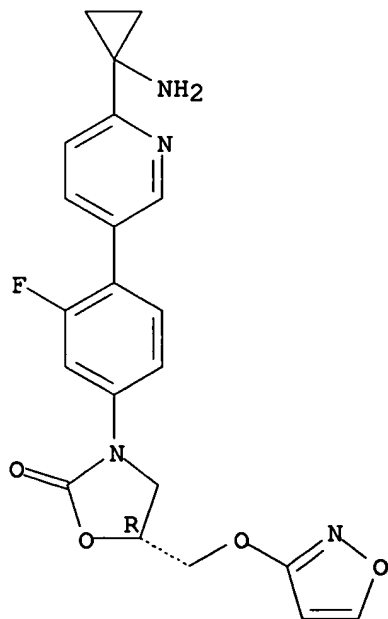
11/539,485



RN 827627-87-2 CAPLUS

CN 2-Oxazolidinone, 3-[4-[6-(1-aminocyclopropyl)-3-pyridinyl]-3-fluorophenyl]-5-[(3-isoxazolyloxy)methyl]-, (5R)- (9CI) (CA INDEX NAME)

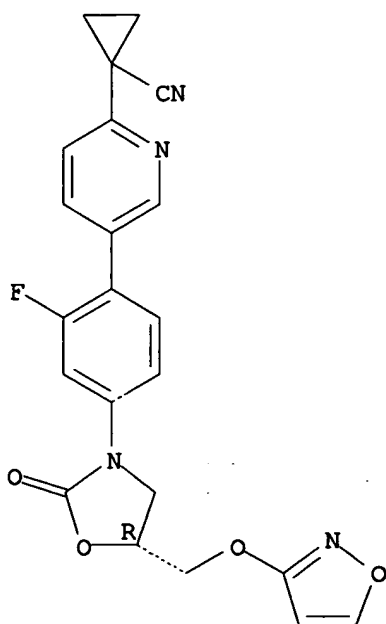
Absolute stereochemistry.



RN 827627-90-7 CAPLUS

CN Cyclopropanecarbonitrile, 1-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RE.CNT 10 THERE ARE 10 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

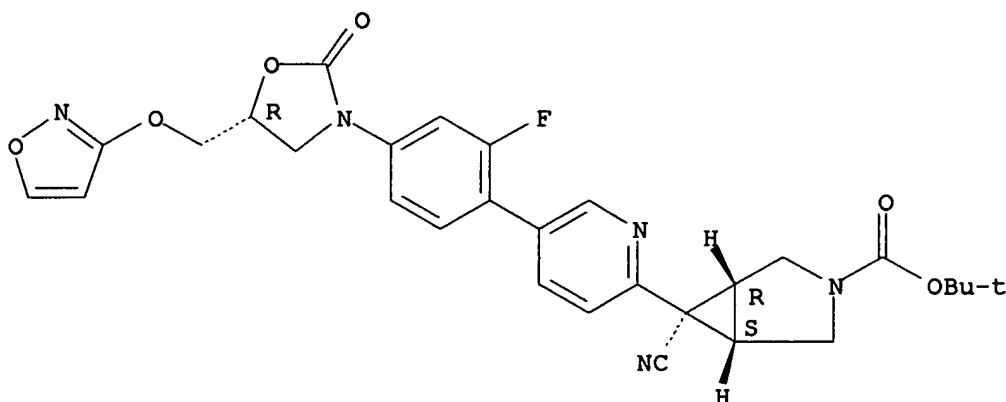
L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN
AN 2005:55212 CAPLUS
DN 142:155937
TI Preparation of cyclopropyl group substituted oxazolidinones as antibiotics
IN Fukuda, Yasumichi
PA Merck & Co., Inc., USA; Kyorin Pharmaceutical Co., Ltd.
SO PCT Int. Appl., 170 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 2

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2005005398	A2	20050120	WO 2004-US20734	20040629
	WO 2005005398	A3	20050428		
	W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
	RW:	BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
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	CA 2530140	AA	20050120	CA 2004-2530140	20040629
	US 2005038092	A1	20050217	US 2004-878637	20040629
	EP 1646629	A2	20060419	EP 2004-777196	20040629
	R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK			
	CN 1816548	A	20060809	CN 2004-80019102	20040629

11/539,485

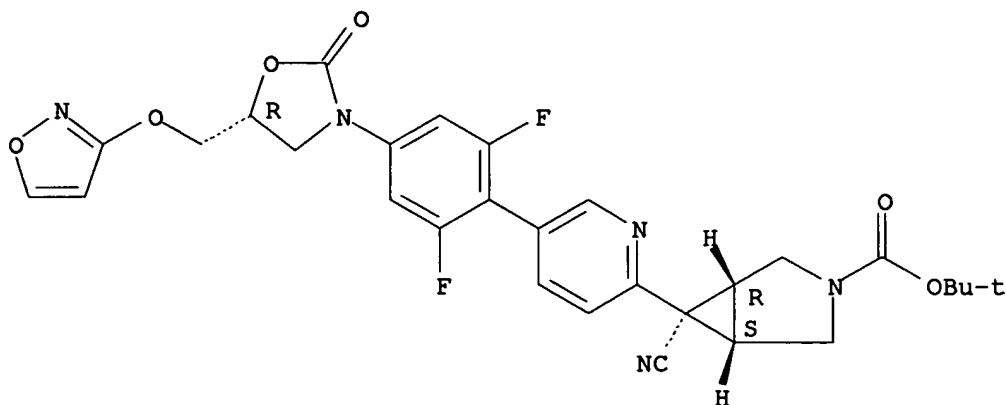
PRAI US 2003-483904P P 20030702
US 2004-546984P P 20040224
WO 2004-US20734 W 20040629
OS CASREACT 142:155937; MARPAT 142:155937
IT 831201-68-4P 831202-25-6P
RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(antibacterial agent; preparation of cyclopropyl-oxazolidinones as antibiotics)
RN 831201-68-4 CAPLUS
CN 3-Azabicyclo[3.1.0]hexane-3-carboxylic acid, 6-cyano-6-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]-, 1,1-dimethylethyl ester, (1 α ,5 α ,6 β)- (9CI)
(CA INDEX NAME)

Absolute stereochemistry.



RN 831202-25-6 CAPLUS
CN 3-Azabicyclo[3.1.0]hexane-3-carboxylic acid, 6-cyano-6-[5-[2,6-difluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]-, 1,1-dimethylethyl ester, (1 α ,5 α ,6 β)- (9CI)
(CA INDEX NAME)

Absolute stereochemistry.



IT 831201-69-5P 831201-72-0P 831202-26-7P
831202-47-2P 831202-76-7P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU

11/539,485

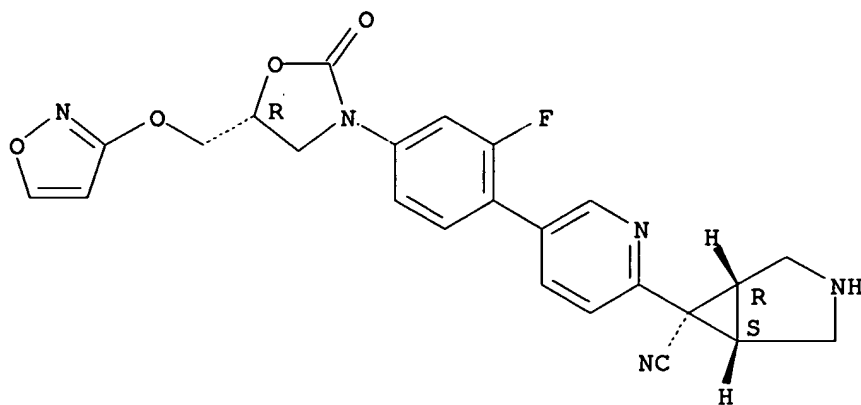
(Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(antibacterial agent; preparation of cyclopropyl-oxazolidinones as antibiotics)

RN 831201-69-5 CAPLUS

CN 3-Azabicyclo[3.1.0]hexane-6-carbonitrile, 6-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]-, monohydrochloride, (1 α ,5 α ,6 β)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

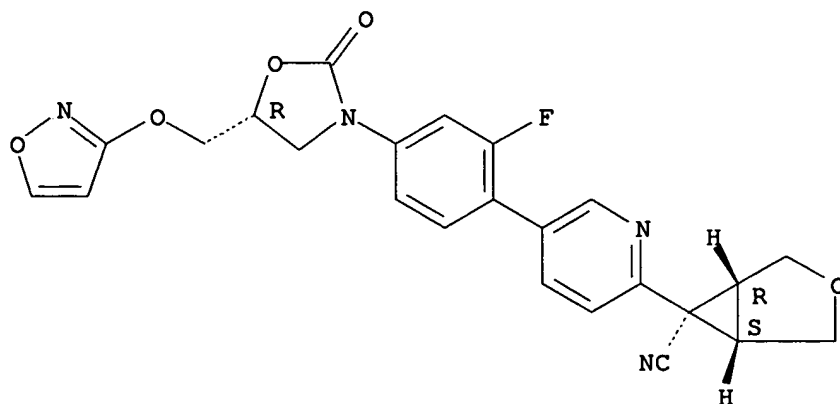


● HCl

RN 831201-72-0 CAPLUS

CN 3-Oxabicyclo[3.1.0]hexane-6-carbonitrile, 6-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]-, (1 α ,5 α ,6 β)- (9CI) (CA INDEX NAME)

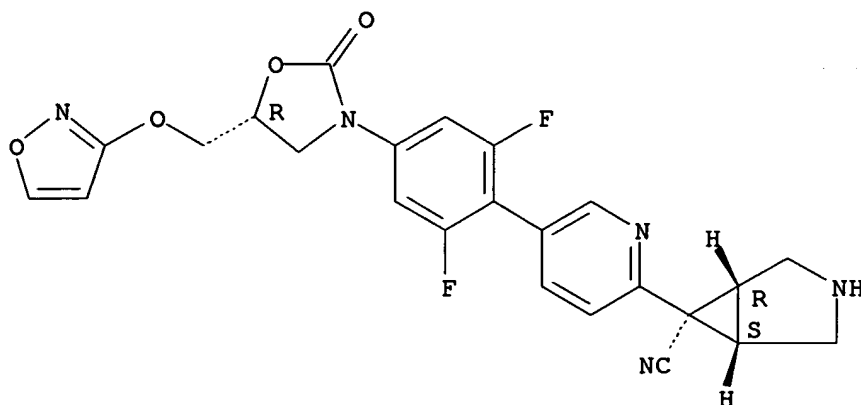
Absolute stereochemistry.



RN 831202-26-7 CAPLUS

CN 3-Azabicyclo[3.1.0]hexane-6-carbonitrile, 6-[5-[2,6-difluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]-, monohydrochloride, (1 α ,5 α ,6 β)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

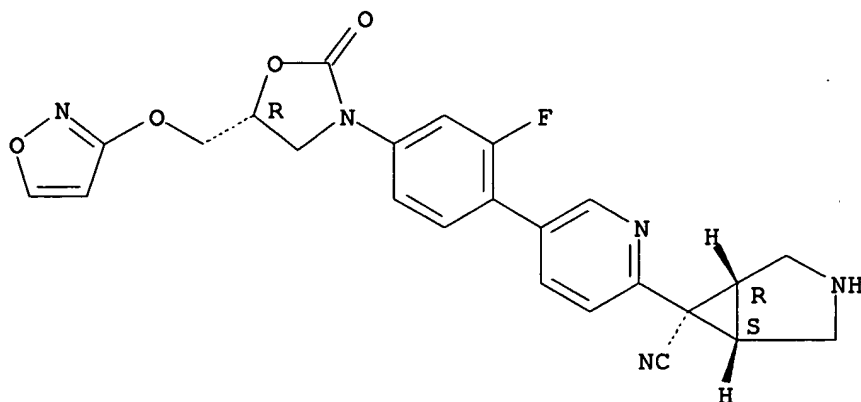


● HCl

RN 831202-47-2 CAPLUS

CN 3-Azabicyclo[3.1.0]hexane-6-carbonitrile, 6-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]-, (1 α ,5 α ,6 β)- (9CI) (CA INDEX NAME)

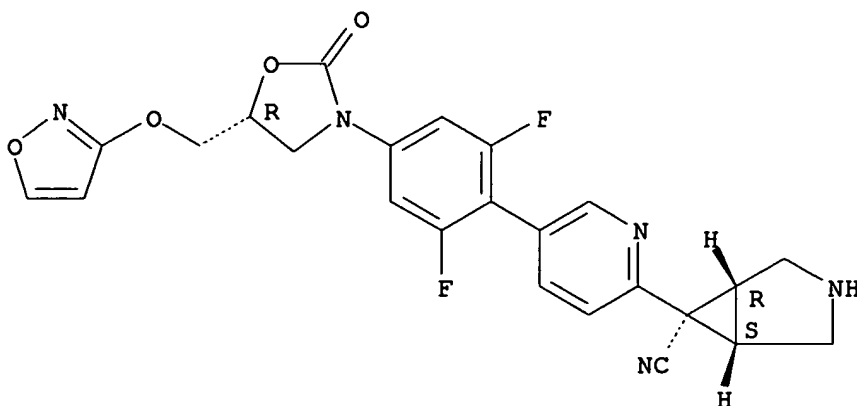
Absolute stereochemistry.



RN 831202-76-7 CAPLUS

CN 3-Azabicyclo[3.1.0]hexane-6-carbonitrile, 6-[5-[2,6-difluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]-, (1 α ,5 α ,6 β)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN
 AN 2004:546503 CAPLUS
 DN 141:89091
 TI A preparation of oxazolidinone derivatives, useful as antibacterial agents
 IN Gravestock, Michael Barry; Hales, Neil James; Turner, Paul
 PA Astrazeneca AB, Swed.; Astrazeneca UK Limited
 SO PCT Int. Appl., 76 pp.
 CODEN: PIXXD2
 DT Patent
 LA English
 FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004056818	A1	20040708	WO 2003-GB5457	20031215
	W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
	RW:	BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
	AU 2003295113	A1	20040714	AU 2003-295113	20031215
	EP 1575945	A1	20050921	EP 2003-786114	20031215
	R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK			
	JP 2006514648	T2	20060511	JP 2004-561619	20031215
	US 2006058317	A1	20060316	US 2005-539485	20050617
PRAI	GB 2002-29518	A	20021219		
	WO 2003-GB5457	W	20031215		

OS MARPAT 141:89091

IT 717116-80-8P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

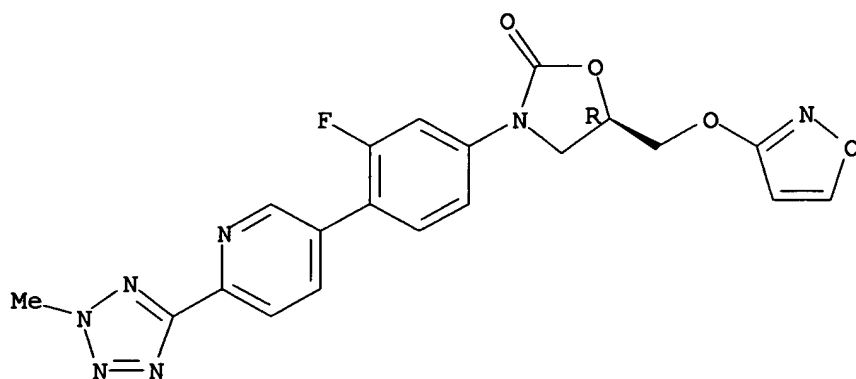
(preparation of oxazolidinone derivs., useful as antibacterial agents)

RN 717116-80-8 CAPLUS

CN 2-Oxazolidinone, 3-[3-fluoro-4-[6-(2-methyl-2H-tetrazol-5-yl)-3-pyridinyl]phenyl]-5-[(3-isoxazolyloxy)methyl]-, (5R)- (9CI) (CA INDEX NAME)

11/539,485

Absolute stereochemistry.



RE.CNT 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

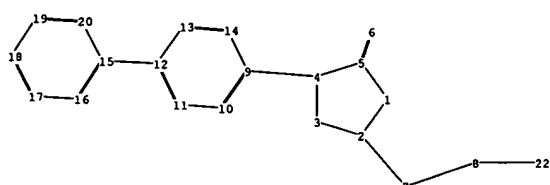
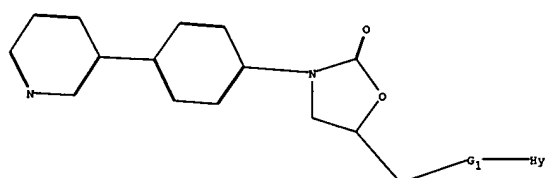
SESSION

FULL ESTIMATED COST

10.99

178.14

STN INTERNATIONAL LOGOFF AT 14:24:16 ON 14 AUG 2006



chain nodes :

6 7 8 22

ring nodes :

1 2 3 4 5 9 10 11 12 13 14 15 16 17 18 19 20

chain bonds :

2-7 4-9 5-6 7-8 8-22 12-15

ring bonds :

1-2 1-5 2-3 3-4 4-5 9-10 9-14 10-11 11-12 12-13 13-14 15-16 15-20 16-17 17-18 18-19 19-20

exact/norm bonds :

1-2 1-5 2-3 3-4 4-5 4-9 5-6 7-8 8-22

exact bonds :

2-7 12-15

normalized bonds :

9-10 9-14 10-11 11-12 12-13 13-14 15-16 15-20 16-17 17-18 18-19 19-20

isolated ring systems :

containing 9 :

G1:O,S

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS7:CLASS8:CLASS9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 22:Atom

Generic attributes :

22:

Saturation

: Unsaturated

11/539,485

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal611bxv

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

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NEWS 6 MAY 11 KOREAPAT updates resume
NEWS 7 MAY 19 Derwent World Patents Index to be reloaded and enhanced
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INPADOC
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and display fields
NEWS 12 JUN 28 Price changes in full-text patent databases EPFULL and PCTFULL
NEWS 13 JUL 11 CHEMSAFE reloaded and enhanced
NEWS 14 JUL 14 FSTA enhanced with Japanese patents
NEWS 15 JUL 19 Coverage of Research Disclosure reinstated in DWPI
NEWS 16 AUG 09 INSPEC enhanced with 1898-1968 archive

NEWS EXPRESS JUNE 30 CURRENT WINDOWS VERSION IS V8.01b, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
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NEWS HOURS STN Operating Hours Plus Help Desk Availability
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NEWS X25 X.25 communication option no longer available

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 14:12:32 ON 14 AUG 2006

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.21

0.21

11/539,485

FILE 'REGISTRY' ENTERED AT 14:12:46 ON 14 AUG 2006
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STRUCTURE FILE UPDATES: 13 AUG 2006 HIGHEST RN 901009-82-3
DICTIONARY FILE UPDATES: 13 AUG 2006 HIGHEST RN 901009-82-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

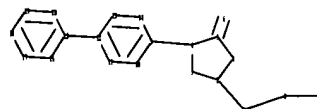
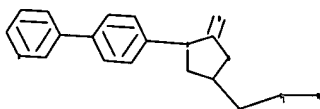
Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10539485.str



chain nodes :

6 7 8 22

ring nodes :

1 2 3 4 5 9 10 11 12 13 14 15 16 17 18 19 20

chain bonds :

2-7 4-9 5-6 7-8 8-22 12-15

ring bonds :

1-2 1-5 2-3 3-4 4-5 9-10 9-14 10-11 11-12 12-13 13-14 15-16 15-20
16-17 17-18 18-19 19-20

11/539,485

exact/norm bonds :

1-2 1-5 2-3 3-4 4-5 4-9 5-6 7-8 8-22

exact bonds :

2-7 12-15

normalized bonds :

9-10 9-14 10-11 11-12 12-13 13-14 15-16 15-20 16-17 17-18 18-19 19-20

isolated ring systems :

containing 9 :

G1:O,S

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:CLASS 9:Atom 10:Atom

11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom

20:Atom 22:Atom

Generic attributes :

22:

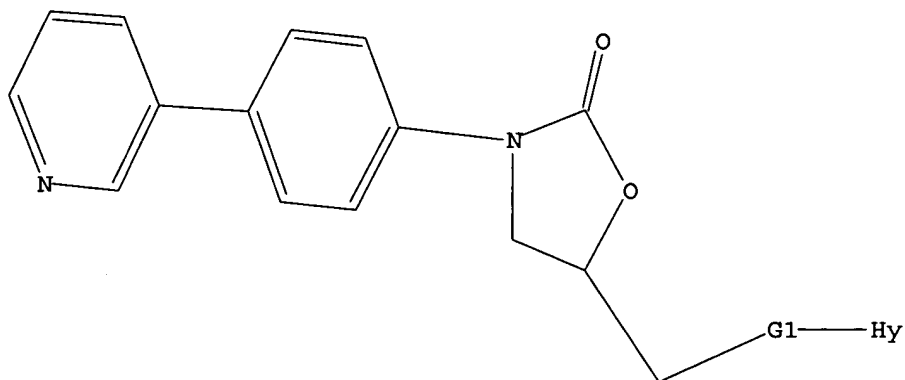
Saturation : Unsaturated

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



G1 O,S

Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss sam

SAMPLE SEARCH INITIATED 14:13:13 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 12 TO ITERATE

100.0% PROCESSED 12 ITERATIONS

SEARCH TIME: 00.00.01

1 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 33 TO 447

PROJECTED ANSWERS: 1 TO 80

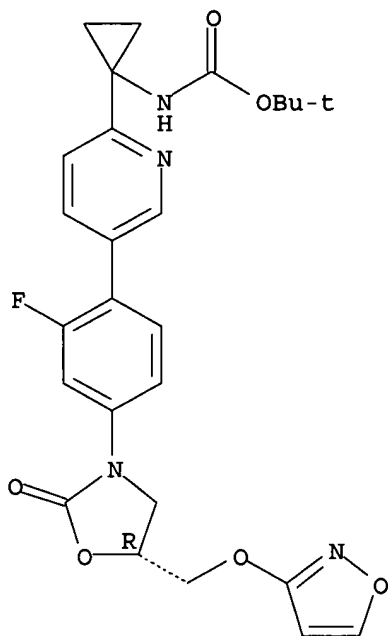
L2 1 SEA SSS SAM L1

11/539,485

=> d scan

L2 1 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Carbamic acid, [1-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]cyclopropyl]-, 1,1-dimethylethyl ester (9CI)
MF C26 H27 F N4 O6

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

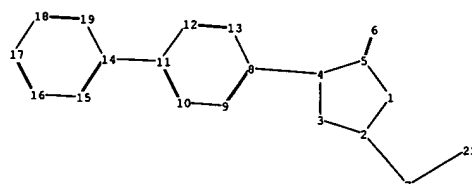
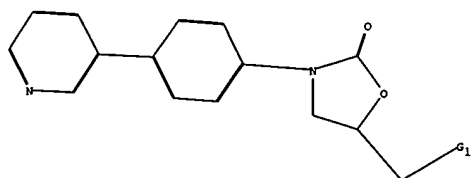
SESSION

FULL ESTIMATED COST

0.44

0.65

STN INTERNATIONAL LOGOFF AT 14:13:31 ON 14 AUG 2006



chain nodes :

6 7 22

ring nodes :

1 2 3 4 5 8 9 10 11 12 13 14 15 16 17 18 19

chain bonds :

2-7 4-8 5-6 7-22 11-14

ring bonds :

1-2 1-5 2-3 3-4 4-5 8-9 8-13 9-10 10-11 11-12 12-13 14-15 14-19 15-16 16-17 17-18 18-19

exact/norm bonds :

1-2 1-5 2-3 3-4 4-5 4-8 5-6 7-22

exact bonds :

2-7 11-14

normalized bonds :

8-9 8-13 9-10 10-11 11-12 12-13 14-15 14-19 15-16 16-17 17-18 18-19

isolated ring systems :

containing 8 :

G1:O,S,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS7:CLASS8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 22:CLASS

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Welcome to STN International! Enter x:x

LOGINID:sssptal611bxv

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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NEWS	12 JUN 28	Price changes in full-text patent databases EPFULL and PCTFULL
NEWS	13 JUL 11	CHEMSAFE reloaded and enhanced
NEWS	14 JUL 14	FSTA enhanced with Japanese patents
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NEWS EXPRESS	JUNE 30	CURRENT WINDOWS VERSION IS V8.01b, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 26 JUNE 2006.
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NEWS X25		X.25 communication option no longer available

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 14:16:28 ON 14 AUG 2006

=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION

11/539,485

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 14:16:37 ON 14 AUG 2006
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STRUCTURE FILE UPDATES: 13 AUG 2006 HIGHEST RN 901009-82-3
DICTIONARY FILE UPDATES: 13 AUG 2006 HIGHEST RN 901009-82-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

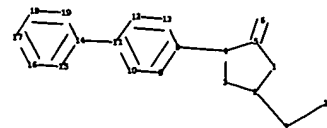
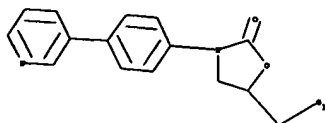
Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10539485-2.str



chain nodes :

11/539,485

6 7 22

ring nodes :

1 2 3 4 5 8 9 10 11 12 13 14 15 16 17 18 19

chain bonds :

2-7 4-8 5-6 7-22 11-14

ring bonds :

1-2 1-5 2-3 3-4 4-5 8-9 8-13 9-10 10-11 11-12 12-13 14-15 14-19 15-16
16-17 17-18 18-19

exact/norm bonds :

1-2 1-5 2-3 3-4 4-5 4-8 5-6 7-22

exact bonds :

2-7 11-14

normalized bonds :

8-9 8-13 9-10 10-11 11-12 12-13 14-15 14-19 15-16 16-17 17-18 18-19

isolated ring systems :

containing 8 :

G1:O,S,N

Match level :

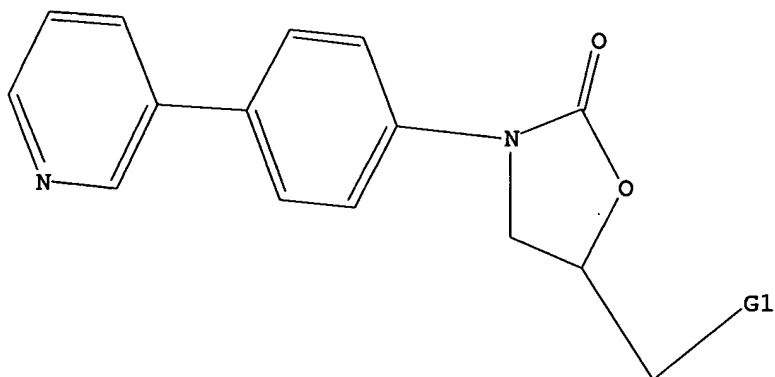
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
22:CLASS

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



G1 O,S,N

Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss sam

SAMPLE SEARCH INITIATED 14:17:00 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 59 TO ITERATE

100.0% PROCESSED 59 ITERATIONS

24 ANSWERS

SEARCH TIME: 00.00.01

11/539,485

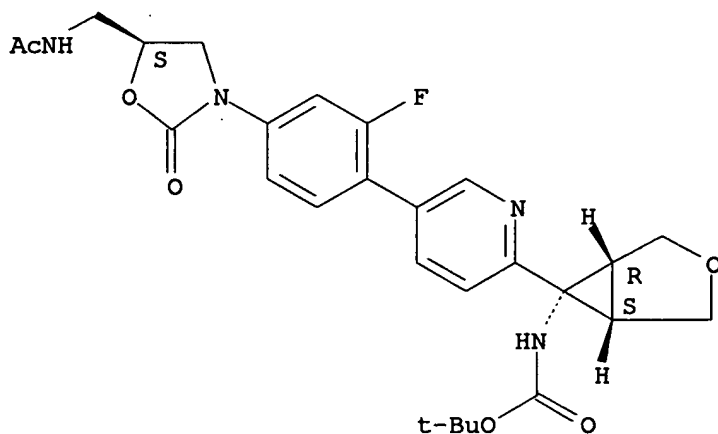
FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 720 TO 1640
PROJECTED ANSWERS: 186 TO 772

L2 24 SEA SSS SAM L1

=> d scan

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Carbamic acid, [(1 α ,5 α ,6 β)-6-[5-[4-[(5S)-5-
[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-fluorophenyl]-2-pyridinyl]-3-
oxabicyclo[3.1.0]hex-6-yl]-, 1,1-dimethylethyl ester (9CI)
MF C27 H31 F N4 O6

Absolute stereochemistry

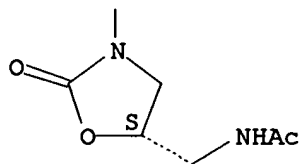
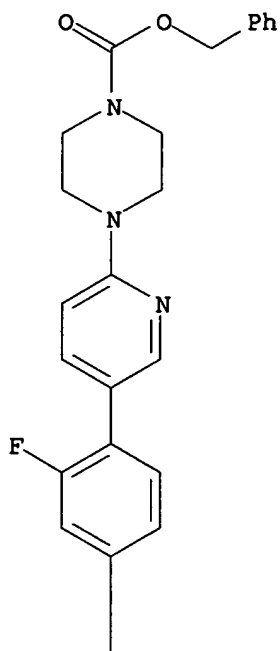


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):23

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN 1-Piperazinecarboxylic acid, 4-[5-[4-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-
oxazolidinyl]-2-fluorophenyl]-2-pyridinyl]-, phenylmethyl ester (9CI)
MF C29 H30 F N5 O5

Absolute stereochemistry.

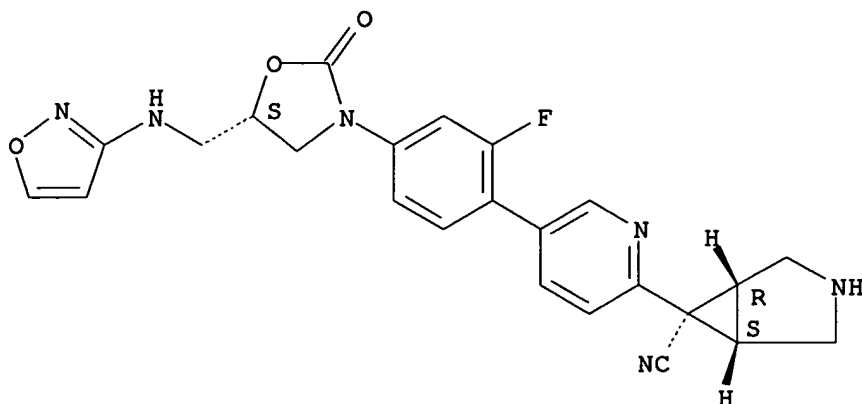


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
 IN 3-Azabicyclo[3.1.0]hexane-6-carbonitrile, 6-[(5S)-5-[2-fluoro-4-[5-[(3-isoxazolylamino)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]-, (1 α ,5 α ,6 β)-(9CI)
 MF C24 H21 F N6 O3
 CI COM

Absolute stereochemistry.

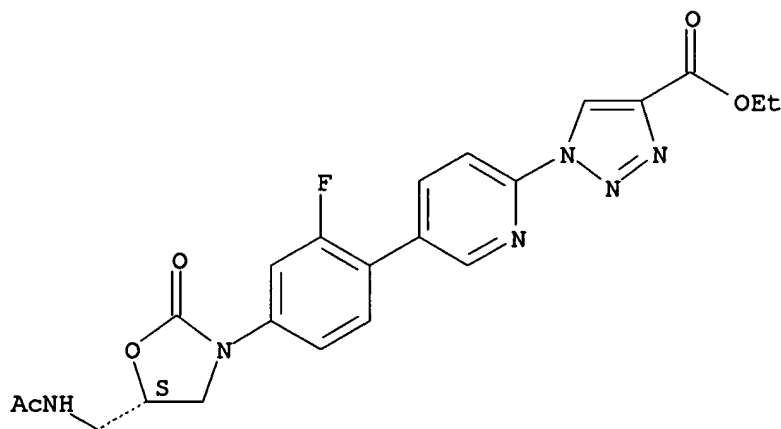
11/539,485



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN 1H-1,2,3-Triazole-4-carboxylic acid, 1-[5-[4-[(5S)-5-[(acetylamino)methyl]-
2-oxo-3-oxazolidinyl]-2-fluorophenyl]-2-pyridinyl]-, ethyl ester (9CI)
MF C22 H21 F N6 O5

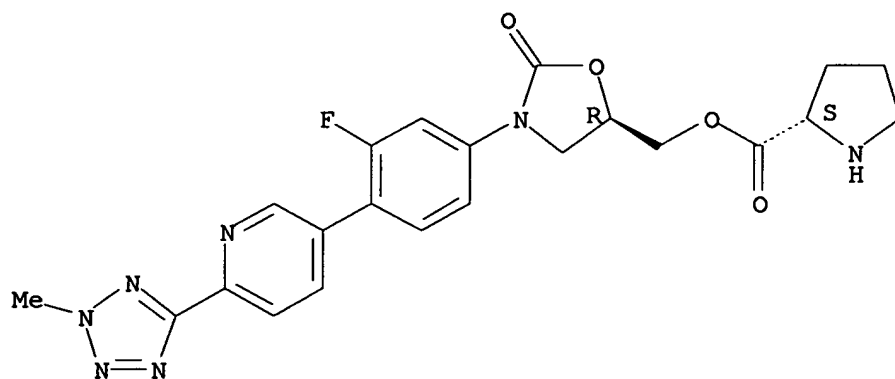
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN L-Proline, [(5R)-3-[3-fluoro-4-[6-(2-methyl-2H-tetrazol-5-yl)-3-
pyridinyl]phenyl]-2-oxo-5-oxazolidinyl]methyl ester, monohydrochloride
(9CI)
MF C22 H22 F N7 O4 . Cl H

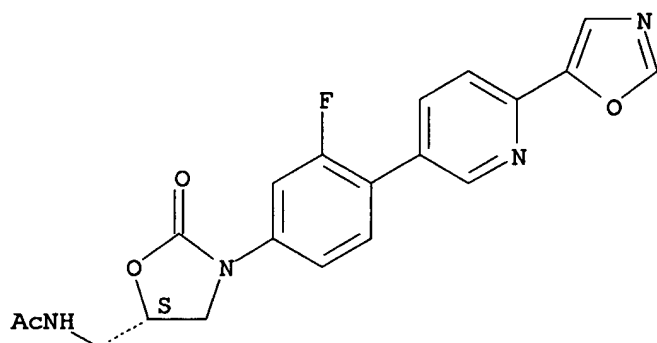
Absolute stereochemistry.



● HCl

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
 IN Acetamide, N-[[[(5S)-3-[3-fluoro-4-[6-(5-oxazolyl)-3-pyridinyl]phenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
 MF C20 H17 F N4 O4

Absolute stereochemistry.

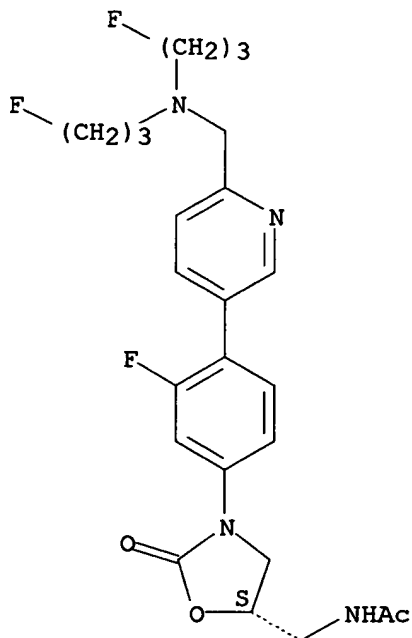


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
 IN Acetamide, N-[[[(5S)-3-[4-[6-[[bis(3-fluoropropyl)amino]methyl]-3-pyridinyl]-3-fluorophenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
 MF C24 H29 F3 N4 O3

Absolute stereochemistry.

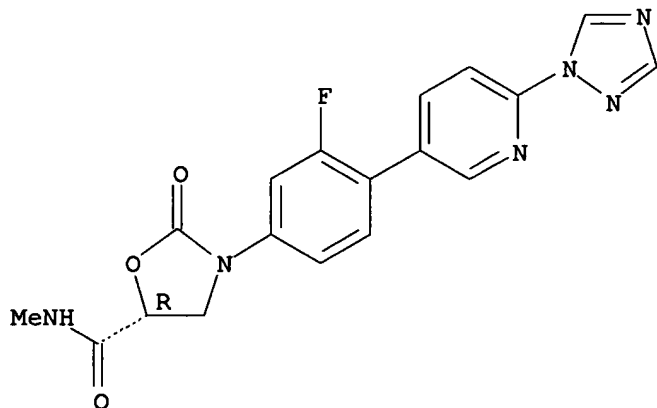
11/539,485



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN 5-Oxazolidinecarboxamide, 3-[3-fluoro-4-[6-(1H-1,2,4-triazol-1-yl)-3-pyridinyl]phenyl]-N-methyl-2-oxo-, (5R)- (9CI)
MF C18 H15 F N6 O3

Absolute stereochemistry.

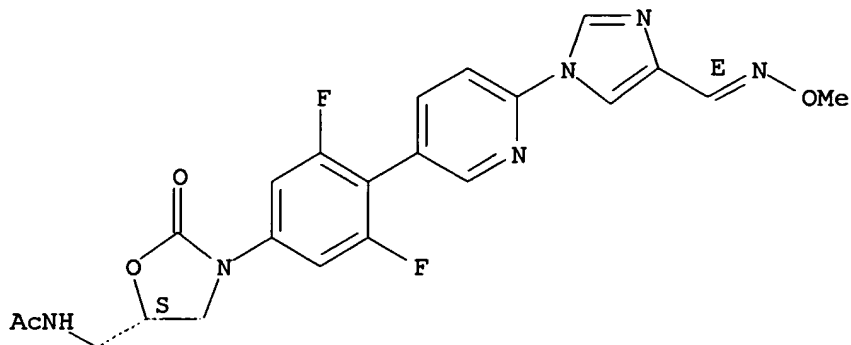


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Acetamide, N-[[(5S)-3-[3,5-difluoro-4-[6-[4-[(E)-(methoxyimino)methyl]-1H-imidazol-1-yl]-3-pyridinyl]phenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
MF C22 H20 F2 N6 O4

11/539,485

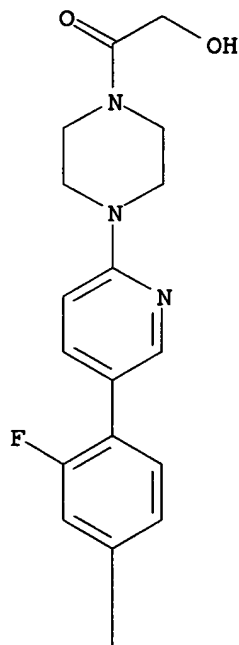
Absolute stereochemistry.
Double bond geometry as shown.

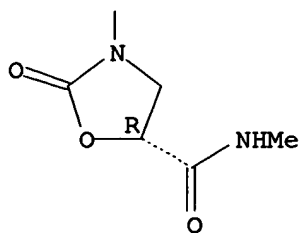


L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN 5-Oxazolidinecarboxamide, 3-[3-fluoro-4-[6-[4-(hydroxyacetyl)-1-piperazinyl]-3-pyridinyl]phenyl]-N-methyl-2-oxo-, (5R)- (9CI)
MF C22 H24 F N5 O5

Absolute stereochemistry.

PAGE 1-A

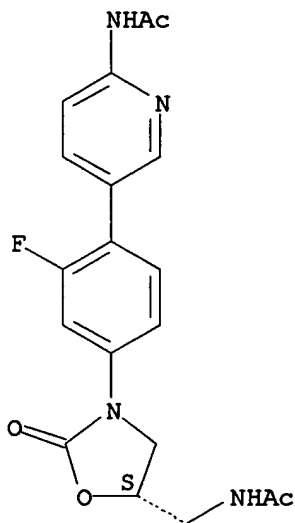




PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
 IN Acetamide, N-[5-[4-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-fluorophenyl]-2-pyridinyl]- (9CI)
 MF C19 H19 F N4 O4

Absolute stereochemistry.

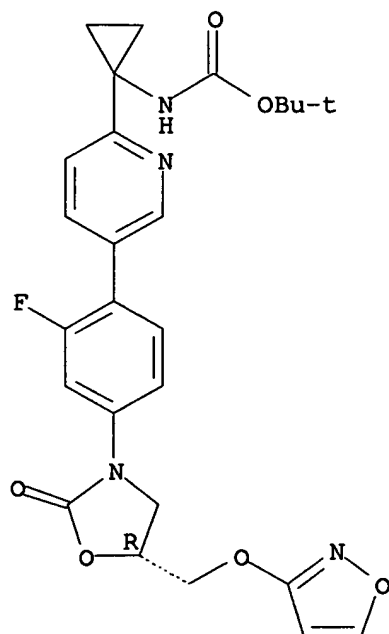


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
 IN Carbamic acid, [1-[5-[2-fluoro-4-[(5R)-5-[(3-isoxazolyloxy)methyl]-2-oxo-3-oxazolidinyl]phenyl]-2-pyridinyl]cyclopropyl]-, 1,1-dimethylethyl ester (9CI)
 MF C26 H27 F N4 O6

Absolute stereochemistry.

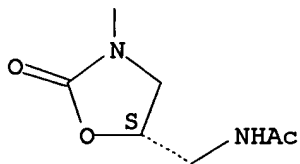
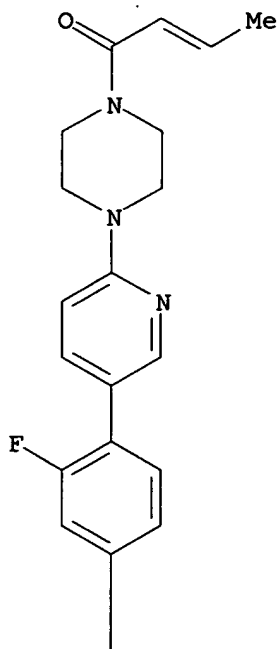
11/539,485



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Acetamide, N-[(5S)-3-[3-fluoro-4-[6-[4-(1-oxo-2-butenyl)-1-piperazinyl]-3-pyridinyl]phenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
MF C25 H28 F N5 O4

Absolute stereochemistry.
Double bond geometry unknown.

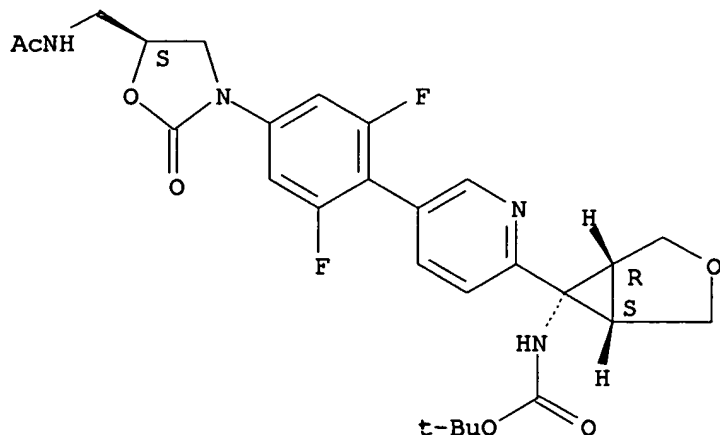


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
 IN Carbamic acid, [(1 α ,5 α ,6 β)-6-[5-[4-[(5S)-5-
 [(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2,6-difluorophenyl]-2-
 pyridinyl]-3-oxabicyclo[3.1.0]hex-6-yl]-, 1,1-dimethylethyl ester (9CI)
 MF C27 H30 F2 N4 O6

Absolute stereochemistry.

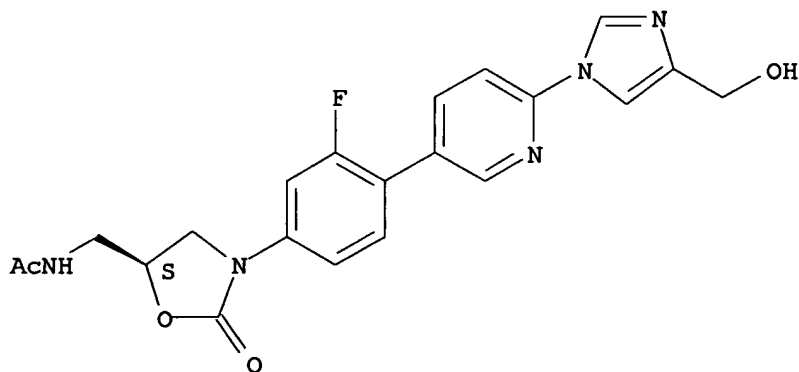
11/539,485



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Acetamide, N-[[(5S)-3-[3-fluoro-4-[6-[4-(hydroxymethyl)-1H-imidazol-1-yl]-3-pyridinyl]phenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
MF C21 H20 F N5 O4

Absolute stereochemistry.

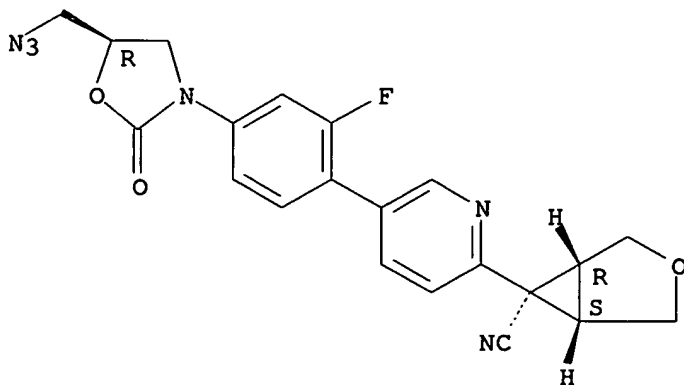


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN 3-Oxabicyclo[3.1.0]hexane-6-carbonitrile, 6-[5-[4-[(5R)-5-(azidomethyl)-2-oxo-3-oxazolidinyl]-2-fluorophenyl]-2-pyridinyl]-, (1 α ,5 α ,6 β)- (9CI)
MF C21 H17 F N6 O3

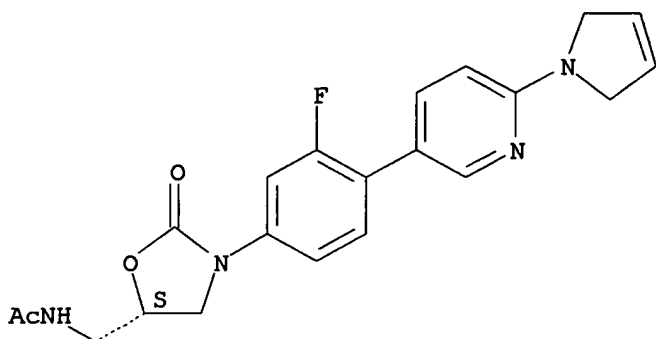
Absolute stereochemistry.

11/539,485



L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Acetamide, N-[[(5S)-3-[4-[6-(2,5-dihydro-1H-pyrrol-1-yl)-3-pyridinyl]-3-fluorophenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
MF C21 H21 F N4 O3

Absolute stereochemistry.

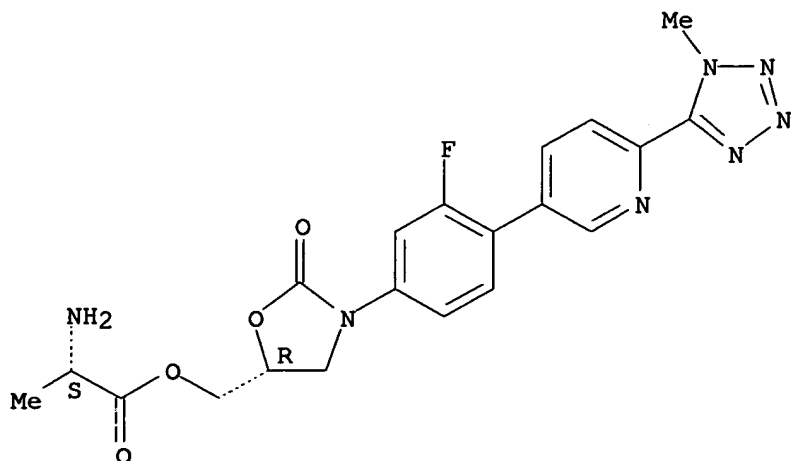


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN L-Alanine, [(5R)-3-[3-fluoro-4-[6-(1-methyl-1H-tetrazol-5-yl)-3-pyridinyl]phenyl]-2-oxo-5-oxazolidinyl]methyl ester (9CI)
MF C20 H20 F N7 O4
CI COM

Absolute stereochemistry.

11/539,485

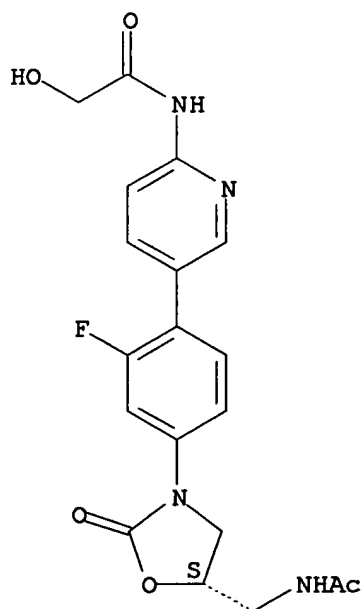


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Cellulose, 2-hydroxypropyl methyl ether, compd. with N-[5-[4-[(5S)-5-
[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-fluorophenyl]-2-pyridinyl]-2-
hydroxyacetamide (9CI)
MF C19 H19 F N4 O5 . x C3 H8 O2 . x C H4 O . x Unspecified

CM 1

Absolute stereochemistry.



CM 2

CM 3

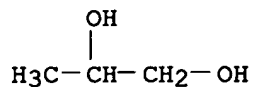
11/539,485

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

CM 4

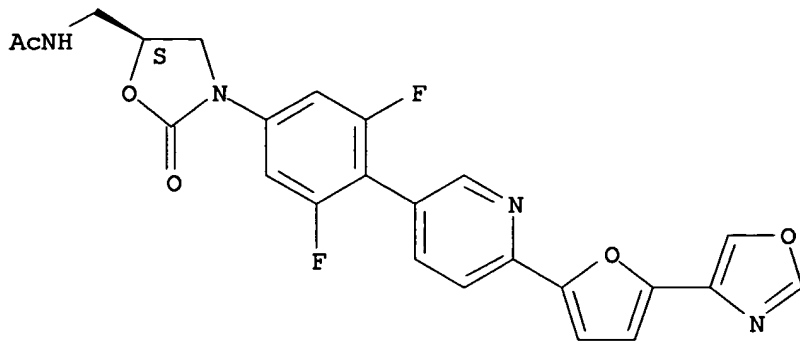
H₃C-OH

CM 5



L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Acetamide, N-[[[(5S)-3-[3,5-difluoro-4-[6-[5-(4-oxazolyl)-2-furanyl]-3-pyridinyl]phenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
MF C24 H18 F2 N4 O5

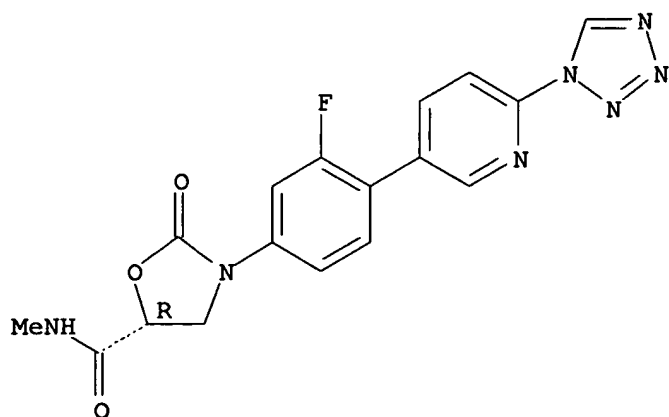
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN 5-Oxazolidinecarboxamide, 3-[3-fluoro-4-[6-(1H-tetrazol-1-yl)-3-pyridinyl]phenyl]-N-methyl-2-oxo-, (5R)- (9CI)
MF C17 H14 F N7 O3

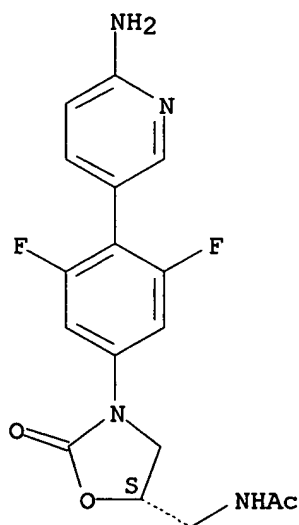
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
 IN Acetamide, N-[[(5S)-3-[4-(6-amino-3-pyridinyl)-3,5-difluorophenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
 MF C17 H16 F2 N4 O3

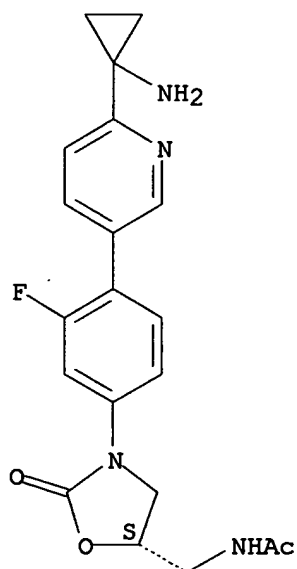
Absolute stereochemistry.



L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
 IN Acetamide, N-[[(5S)-3-[4-[6-(1-aminocyclopropyl)-3-pyridinyl]-3-fluorophenyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI)
 MF C20 H21 F N4 O3

Absolute stereochemistry.

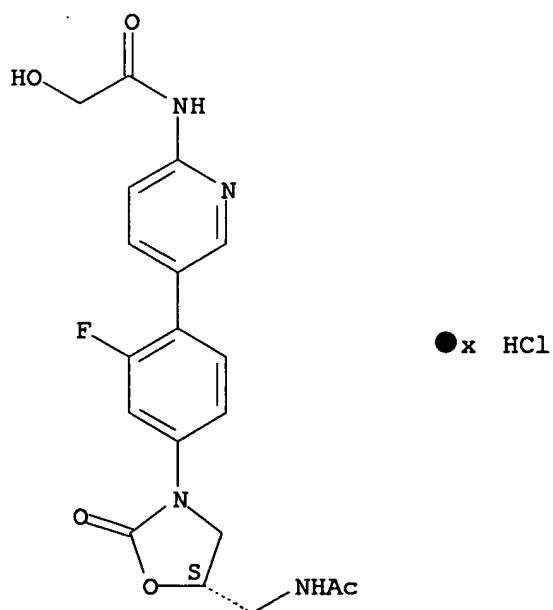
11/539,485



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 24 ANSWERS REGISTRY COPYRIGHT 2006 ACS on STN
IN Acetamide, N-[5-[4-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-fluorophenyl]-2-pyridinyl]-2-hydroxy-, hydrochloride (9CI)
MF C19 H19 F N4 O5 . x Cl H

Absolute stereochemistry.



ALL ANSWERS HAVE BEEN SCANNED

11/539,485

=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

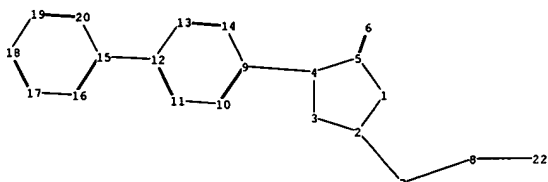
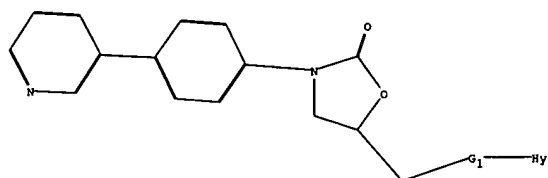
SESSION

FULL ESTIMATED COST

0.88

1.09

STN INTERNATIONAL LOGOFF AT 14:17:33 ON 14 AUG 2006



chain nodes :

6 7 8 22

ring nodes :

1 2 3 4 5 9 10 11 12 13 14 15 16 17 18 19 20

chain bonds :

2-7 4-9 5-6 7-8 8-22 12-15

ring bonds :

1-2 1-5 2-3 3-4 4-5 9-10 9-14 10-11 11-12 12-13 13-14 15-16 15-20 16-17 17-18 18-19 19-20

exact/norm bonds :

1-2 1-5 2-3 3-4 4-5 4-9 5-6 7-8 8-22

exact bonds :

2-7 12-15

normalized bonds :

9-10 9-14 10-11 11-12 12-13 13-14 15-16 15-20 16-17 17-18 18-19 19-20

isolated ring systems :

containing 9 :

G1:O,S

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS7:CLASS8:CLASS9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 22:Atom

Generic attributes :

22:

Saturation

: Unsaturated